

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120320\
 Data File : BN012849.D
 Acq On : 03 Dec 2020 16:32
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.411

Quant Time: Dec 03 17:01:32 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 03 15:27:55 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	505	0.40	ng/ul	0.00
2) Naphthalene-d8	10.17	136	1895	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1101	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.82	188	2411	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	2056	0.40	ng/ul	0.00
20) Perylene-d12	23.15	264	2017	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.78	152	1135	0.40	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	2594	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.21	128	2185	0.398	ng/ul	99
5) 2-Methylnaphthalene	11.85	142	1413	0.392	ng/ul	100
7) Acenaphthylene	13.78	152	2067	0.397	ng/ul	99
8) Acenaphthene	14.13	153	1720	0.402	ng/ul	98
9) Fluorene	15.12	166	1966	0.401	ng/ul	99
11) Pentachlorophenol	16.49	266	251	0.359	ng/ul	97
12) Phenanthrene	16.87	178	3169	0.392	ng/ul	98
13) Anthracene	16.96	178	2784	0.405	ng/ul	99
15) Fluoranthene	18.90	202	3734	0.389	ng/ul	97
17) Pyrene	19.26	202	3821	0.393	ng/ul	99
18) Benzo(a)anthracene	21.03	228	3317	0.390	ng/ul	99
19) Chrysene	21.08	228	3894	0.414	ng/ul	99
21) Benzo(b)fluoranthene	22.53	252	3664	0.393	ng/ul	99
22) Benzo(k)fluoranthene	22.57	252	4165	0.404	ng/ul	99
23) Benzo(a)pyrene	23.06	252	3356	0.388	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.22	276	4583	0.403	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.23	278	3622	0.403	ng/ul	98
26) Benzo(g,h,i)perylene	25.84	276	4106	0.399	ng/ul	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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